

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121418\
 Data File : BF111434.D
 Acq On : 14 Dec 2018 22:20
 Operator : JU/SJ
 Sample : J6371-02 2X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EB07-1-2

Manual Integrations
 APPROVED

Sohil
 12/17/2018 2:36:57 PM

Quant Time: Dec 14 22:52:17 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 14 13:26:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	185037	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	719268	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	285630	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	426202	20.00	ng	0.00
76) Chrysene-d12	13.95	240	342761	20.00	ng	0.00
87) Perylene-d12	15.39	264	247414	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	209446	18.84	ng	0.00
7) Phenol-d6	6.43	99	705989	47.73	ng	0.00
23) Nitrobenzene-d5	7.35	82	462731	38.31	ng	0.00
42) 2,4,6-Tribromophenol	10.62	330	23198	9.07	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	789858	42.71	ng	0.00
79) Terphenyl-d14	12.90	244	556828	38.15	ng	0.00
Target Compounds						
10) Phenol	6.45	94	50027	3.036	ng	93
50) Dimethylphthalate	9.55	163	60925	2.962	ng	# 99
71) Phenanthrene	11.34	178	349463	15.913	ng	93
72) Anthracene	11.39	178	95286	4.243	ng	93
75) Fluoranthene	12.53	202	750724	34.942	ng	97
78) Pyrene	12.76	202	697582	28.866	ng	95
81) Benzo(a)anthracene	13.94	228	365142	19.592	ng	99
83) Chrysene	13.97	228	338119	17.222	ng	96
86) Indeno(1,2,3-cd)pyrene	16.81	276	129750m	9.157	ng	
88) Benzo(b)fluoranthene	14.98	252	367561m	25.482	ng	
89) Benzo(k)fluoranthene	15.00	252	94756m	6.875	ng	
90) Benzo(a)pyrene	15.33	252	235204	18.082	ng	98
91) Dibenzo(a,h)anthracene	16.82	278	31597	3.079	ng	# 85
92) Benzo(g,h,i)perylene	17.23	276	131013	13.008	ng	93

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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